Zadatak 1: Zaključivanje u statičkoj Bejzovoj mreži

Za Bejzovu mrežu datu u postavci domaćeg, potrebno je odrediti uslovnu verovatnoću $p(c^+|f^+,g^+)$.

Ovo je potrebno uraditi sledećim metodama:

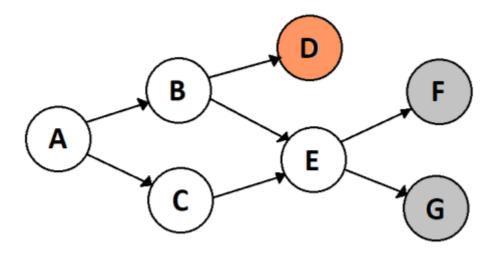
- · metodom eliminacije,
- · uzorkovanjem sa odbacivanjem,
- Gibbs-ovim uzorkovanjem (neparan indeks).

1.1. Prikaz grafa Bejzove mreže

```
import matplotlib.pyplot as plt
import matplotlib.image as mpimg

img = mpimg.imread('./BM_model.png')

plt.imshow(img); plt.axis('off'); plt.show()
```



1.2. Procena $p(c^+|f^+,g^+)$ metodom eliminacije

Metoda eliminacije je najjednostavnija metoda tačnog zaključivanja u Bejzovim mrežama. Ne uvodi aproksimacije, već izračunava tačne verovatnoće događaja koristeći osobine uslovne verovatnoće i strukturu date mreže.

Da bismo našli datu uslovnu raspodelu, potrebno je primeniti Bejzovo pravilo:

$$p(C|f^+,g^+) \propto p(C,f+,g+) = \sum\limits_{a,b,e} p(a) \cdot p(b|a) \cdot p(C|a) \cdot p(e|b,C) \cdot p(f^+|e) \cdot p(g^+|e)$$

U prethodnom izrazu je potrebno **marginalizovati** promenljive a, b i e. Ovo možemo uraditi pomoću više redosleda biranja promenljivih. Pokazuje se da će redosled A, E, B da ima manje potrebnih računskih operacija, pa ćemo tim redosledom ih marginalizovati.

• Eliminisanje promenljive a:

$$\phi_7(a,b,c) = p(a) \cdot p(b|a) \cdot p(C|a)$$
 $f_7(b,c) = \sum\limits_a \phi_7(a,b,c)$

b	С	а	$\phi_7(b,c,a)$	$f_7(b,c)$
0	0	0	0.6 * 0.7 * 0.7 = 0.294	0.318
0	0	1	0.4 * 0.2 * 0.3 = 0.024	
0	1	0	0.6 * 0.7 * 0.3 = 0.126	0.182
0	1	1	0.4 * 0.2 * 0.7 = 0.056	
1	0	0	0.6 * 0.3 * 0.7 = 0.126	0.222
1	0	1	0.4 * 0.8 * 0.3 = 0.096	
1	1	0	0.6 * 0.3 * 0.3 = 0.054	0.278
1	1	1	0.4 * 0.8 * 0.7 = 0.224	

$$p(C,f+,g+) = \sum\limits_{b,e} f_7(b,c) \cdot p(e|b,C) \cdot p(f^+|e) \cdot p(g^+|e)$$

• Eliminisanje promenljive e:

$$egin{aligned} \phi_8(b,c,e) &= p(e|b,c) \cdot p(f^+|e) \cdot p(g^+|e) \ f_8(b,c) &= \sum\limits_e \phi_8(b,c,e) \end{aligned}$$

b
 c
 e

$$\phi_8(b,c,e)$$
 $f_8(b,c)$

 0
 0
 0.1 * 0.6 * 0.2 = 0.012
 0.264

 0
 0
 1
 0.9 * 0.4 * 0.7 = 0.252

 0
 1
 0
 0.5 * 0.6 * 0.2 = 0.06
 0.2

 0
 1
 1
 0.5 * 0.4 * 0.7 = 0.14
 0.248

 1
 0
 0
 0.2 * 0.6 * 0.2 = 0.024
 0.248

 1
 0
 1
 0.8 * 0.4 * 0.7 = 0.224
 0.1552

 1
 1
 0.2 * 0.4 * 0.7 = 0.056
 0.056

$$p(C,f+,g+) = \sum\limits_b f_7(b,c) \cdot f_8(b,c)$$

• Eliminisanje promenljive b:

$$\phi_9(b,c) = f_7(b,c) \cdot f_8(b,c)$$

$$f_9(c) = \sum\limits_b \phi_9(b,c)$$

b	c	$\phi_9(b,c)$			
0	0	0.318 * 0.264 = 0.083952			
0	1	0.182 * 0.2 = 0.0364			
1	0	0.222 * 0.248 = 0.055056			
1	1	0.278 * 0.152 = 0.042256			

$$p(c^+|f^+,g^+)=lpha\cdot(0.0364+0.042256)=lpha\cdot0.078656$$
 $p(c^-|f^+,g^+)=lpha\cdot(0.083952+0.055056)=lpha\cdot0.139008$ $p(c^+|f^+,g^+)+p(c^-|f^+,g^+)=lpha\cdot0.217664=1$ $p(c^+|f^+,g^+)pprox0.361364$

1.3. Procena $p(c^+|f^+,g^+)$ uzorkovanjem sa odbacivanjem

Prethodna metoda je bila 'tačna metoda' za određivanje verovatnoće. Naredne dve metode predstavljaju Monte Karlo metode procene i zasnivaju se na uzorkovanju promenljivih.

Najpre cemo uzorkovati sa odbacivanjem. Ova metoda podrazumeva da na osnovu datih raspodela u postavci domaćeg generišemo odbirke promenljivih a zatim iz takvog uzorka procenimo verovatnoću kao:

$$p(c^+|f^+,g^+) pprox rac{\#(*,*,c^+,*,*,f^+,g^+)}{\#(*,*,*,*,*,f^+,g^+)}$$

Odbacivanje

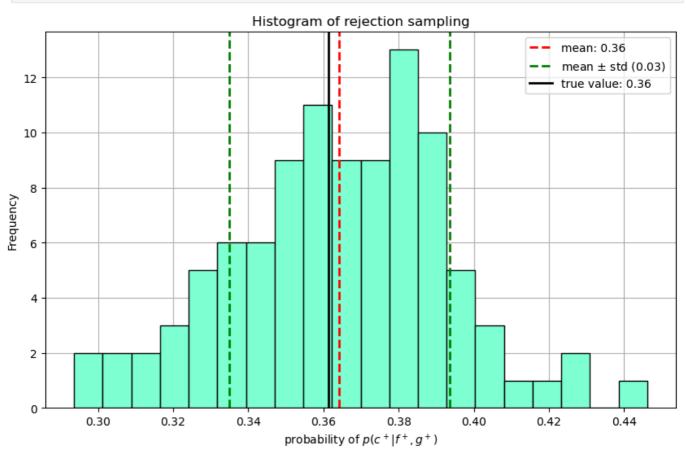
Prilikom uzorkovanja, ukoliko bismo išli po promenljivama abecednim redom, kada dobijemo kao uzorak za $f=f^-$, onda nema potrebe da dalje uzorkujemo i promenljivu g. Ova ušteda se naziva odbacivanje i mi ćemo sa ovim poboljšanjem da realizujemo procenu.

```
In [3]: from random import random
        import numpy as np
        # Probability tables
        table a = np.array([0.6])
        table_b = np.array([0.7, 0.2])
        table_c = np.array([0.7, 0.3])
        table_d = np.array([0.4, 0.8])
        table_e = np.array([0.1, 0.5, 0.2, 0.8])
        table_f = np.array([0.4, 0.6])
        table_g = np.array([0.8, 0.3])
        # Define functions for generating an output based on inputs
        def generate_A():
            r = random()
            if r < 0.6:
                return False
            return True
```

```
def generate_B(a: bool):
              r = random()
              compare_value = table_b[np.packbits([a], bitorder='little')]
              if r<compare_value:</pre>
                  return False
              return True
          def generate_C(a: bool):
             r = random()
              compare_value = table_c[np.packbits([a], bitorder='little')]
              if r<compare_value:</pre>
                  return False
              return True
          def generate_D(b: bool):
             r = random()
              compare_value = table_d[np.packbits([b], bitorder='little')]
              if r<compare value:</pre>
                  return False
              return True
          def generate_E(b: bool, c:bool):
             r = random()
              compare_value = table_e[np.packbits([c, b], bitorder='little')]
              if r<compare value:</pre>
                  return False
              return True
          def generate_F(e: bool):
             r = random()
              compare_value = table_f[np.packbits([e], bitorder='little')]
              if r<compare_value:</pre>
                  return False
              return True
          def generate_G(e: bool):
              r = random()
              compare_value = table_g[np.packbits([e], bitorder='little')]
              if r<compare value:</pre>
                  return False
              return True
In [35]: N_rejection = 1200
```

```
Nr = 100
rejection_estimations = np.empty(Nr, dtype=float)
for i in range(Nr):
    sample count = 0
    specific_count = 0
    for j in range(N_rejection):
        a = generate_A()
        b = generate_B(a)
        c = generate_C(a)
        e = generate_E(b, c)
        f = generate_F(e)
        if not f:
            continue
        g = generate_G(e)
        if f and g:
            sample_count+=1
            if c:
                specific_count+=1
```

```
In [36]:
         import matplotlib.pyplot as plt
         mean_rejection = np.mean(rejection_estimations)
         std_dev_rejection = np.std(rejection_estimations)
         x = 0.361364
         plt.figure(figsize=(10,6))
         plt.rc('axes', axisbelow=True)
         plt.hist(rejection_estimations, color='aquamarine', bins=20, edgecolor='black')
         plt.axvline(mean_rejection, color='red', linestyle='dashed', linewidth=2, label=f'mean: {mean_
         plt.axvline(mean_rejection - std_dev_rejection, color='green', linestyle='dashed', linewidth=
         plt.axvline(mean_rejection + std_dev_rejection, color='green', linestyle='dashed', linewidth=
         plt.axvline(x, color='black', linestyle='solid', linewidth=2, label=f'true value: {x:.2f}')
         plt.legend()
         plt.xlabel('probability of $p(c^+|f^+,g^+)$')
         plt.ylabel('Frequency')
         plt.title('Histogram of rejection sampling')
         plt.grid(True)
         plt.show()
```



1.4. Procena $p(c^+|f^+,g^+)$ Gibbs-ovim uzorkovanjem

Gibbs-ovo uzorkovanje je metoda Monte Karlo uzorkovanja. Predstavlja posebnu vrstu MCMC-a (Markov Chain Monte Carlo). Metoda ima za zadatak da uzorkuje iz raspodele $p(X)=p(x_1,x_2,\ldots,x_n)$. Princip rada je sledeći:

```
1. inicijalizacija: X^{(0)}=(x_1^{(0)},x_2^{(0)},\dots,x_n^{(0)})
2. iteracije i=1,\dots,N :
```

$$egin{aligned} x_1^{(i)} &\sim p(x_1|x_2^{(i-1)},x_3^{(i-1)},\dots,x_n^{(i-1)}) \ &x_2^{(i)} &\sim p(x_2|x_1^{(i)},x_3^{(i-1)},\dots,x_n^{(i-1)}) \ & \cdots \ &x_n^{(i)} &\sim p(x_n|x_1^{(i)},x_2^{(i)},\dots,x_{n-1}^{(i-1)}) \end{aligned}$$

Primenom ove metode ćemo uzorkovati uzorak za našu Bejzovu mrežu. Da bismo ovo postigli, potrebno je da nađemo uslovne raspodele svih promenljivih mreže pri datim svim ostalim promenljivama. Za ove potrebe ćemo koristiti sledeću osobinu:

$$MP(X_i) = R \cup D \cup S$$
 $p(X_i|x_{j \neq i}) \propto p(D|X_i,S) \cdot p(X_i|R)$

1. Raspodela $p(A|b,c,d,e,f^+,g^+)$:

$$p(A|...) \propto p(b,c|A) \cdot p(A)$$

b
 c

$$p(a^+ \parallel ...)$$
 $p(a^- \parallel ...)$
 α^{-1}
 $p(a^+ \parallel b, c)$

 0
 0
 0.024 * α
 0.294 * α
 0.318
 0.0755

 0
 1
 0.056 * α
 0.126 * α
 0.182
 0.3077

 1
 0
 0.096 * α
 0.126 * α
 0.222
 0.4324

 1
 1
 0.224 * α
 0.054 * α
 0.278
 0.8058

2. Raspodela $p(B|a,c,d,e,f^+,g^+)$:

$$p(B|...) \propto p(e|B,c) \cdot p(d|B) \cdot p(B|a)$$

а	c	d	е	$p(b^+\ \dots)/lpha$	$p(b^-\ \dots)/lpha$	$lpha^{-1}$	$p(b^+ \ a, c, d, e)$
0	0	0	0	0.048	0.028	0.076	0.6316
0	0	0	1	0.192	0.252	0.444	0.4324
0	0	1	0	0.012	0.042	0.054	0.2222
0	0	1	1	0.048	0.378	0.426	0.1127
0	1	0	0	0.192	0.140	0.332	0.5783
0	1	0	1	0.048	0.140	0.188	0.2553
0	1	1	0	0.048	0.210	0.258	0.1860
0	1	1	1	0.012	0.210	0.222	0.0540
1	0	0	0	0.128	0.008	0.136	0.9412
1	0	0	1	0.512	0.072	0.584	0.8767
1	0	1	0	0.032	0.012	0.044	0.7272
1	0	1	1	0.128	0.108	0.236	0.5424
1	1	0	0	0.512	0.040	0.552	0.9275

а	c	d	e	$p(b^+\ \dots)/lpha$	$p(b^-\ \dots)/lpha$	$lpha^{-1}$	$p(b^+ \ a, c, d, e)$

1	1	0	1	0.128	0.040	0.168	0.7619
1	1	1	0	0.128	0.060	0.188	0.6808
1	1	1	1	0.032	0.060	0.092	0.3478

3. Raspodela $p(C|a,b,d,e,f^+,g^+)$:

$$p(C|...) \propto p(e|b,C) \cdot p(C|a)$$

a
 b
 e

$$p(c^+||...)/\alpha$$
 $p(c^-||...)/\alpha$
 $p(c^-||...)/\alpha$
 $p(c^+||a,b,e)$

 0
 0
 0
 0.15
 0.07
 0.22
 0.6818

 0
 0
 1
 0.15
 0.63
 0.78
 0.1923

 0
 1
 0
 0.24
 0.14
 0.38
 0.6316

 0
 1
 1
 0.06
 0.56
 0.62
 0.0968

 1
 0
 0
 0.35
 0.03
 0.38
 0.9210

 1
 0
 1
 0.35
 0.27
 0.62
 0.5645

 1
 1
 0
 0.56
 0.06
 0.62
 0.9032

 1
 1
 1
 0.14
 0.24
 0.38
 0.3684

- 4. Raspodela $p(D|a,b,c,e,f^+,g^+)=p(D|b)$
- 5. Raspodela $p(E|a,b,c,d,f^+,g^+)$:

$$p(E|\dots) \propto p(f^+|E) \cdot p(g^+|E) \cdot p(E|b,c)$$

b
 c

$$p(e^+ \parallel ...)/\alpha$$
 $p(e^- \parallel ...)/\alpha$
 α^{-1}
 $p(e^+ \parallel b, c)$

 0
 0
 0.252
 0.012
 0.264
 0.9545

 0
 1
 0.140
 0.060
 0.200
 0.7000

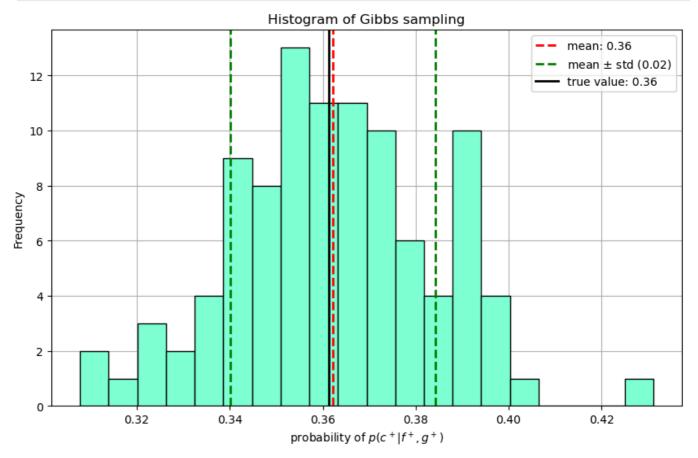
 1
 0
 0.224
 0.024
 0.248
 0.9032

 1
 1
 0.056
 0.096
 0.152
 0.3684

```
gibbs_table_e = np.array([0.9545, 0.7, 0.9032, 0.3684])
def gibbs_generate_A(b:bool, c:bool):
   r = random()
    compare_value = gibbs_table_a[np.packbits([c, b], bitorder='little')]
    if r<compare_value:</pre>
        return True
    return False
def gibbs_generate_B(a:bool, c:bool, d:bool, e:bool):
    r = random()
    compare_value = gibbs_table_b[np.packbits([e, d, c, a], bitorder='little')]
    if r<compare_value:</pre>
        return True
    return False
def gibbs_generate_C(a:bool, b:bool, e:bool):
    r = random()
    compare_value = gibbs_table_c[np.packbits([e, b, a], bitorder='little')]
    if r<compare value:</pre>
        return True
    return False
def gibbs_generate_D(b: bool):
   return generate_D(b)
def gibbs_generate_E(b:bool, c:bool):
    r = random()
   compare_value = gibbs_table_e[np.packbits([c, b], bitorder='little')]
   if r<compare_value:</pre>
        return True
    return False
```

```
In [18]: # Gibbs sampling
         from random import choice
         Nr = 100
         N_gibbs = 1000
         burnout_percentage = 10 # percent of samples we reject from the beginning of gibbs sampling
         burnout rejection count = N gibbs*burnout percentage/100
         gibbs_estimations = np.empty(Nr, dtype=float)
         for i in range(Nr):
             # Initializing
             a, b, c, d, e = [choice([True, False]) for _ in range(5)]
             # Iterations
             count = 0
             for j in range(1,N_gibbs):
                 # Generating a point
                 a = gibbs_generate_A(b, c)
                 b = gibbs_generate_B(a,c,d,e)
                 c = gibbs_generate_C(a,b,e)
                 d = gibbs_generate_D(b)
                 e = gibbs_generate_E(b,c)
                 if c and j>=burnout_rejection_count:
                     count+=1
             # Finding the probability(p(c+|f+, g+))
             gibbs_estimations[i] = count/(N_gibbs-burnout_rejection_count)
```

```
In [19]: import matplotlib.pyplot as plt
         mean_gibbs = np.mean(gibbs_estimations)
         std_dev_gibbs = np.std(gibbs_estimations)
         x = 0.361364
         plt.figure(figsize=(10,6))
         plt.rc('axes', axisbelow=True)
         plt.hist(gibbs_estimations, color='aquamarine', bins=20, edgecolor='black')
         plt.axvline(mean_gibbs, color='red', linestyle='dashed', linewidth=2, label=f'mean: {mean_gib|
         plt.axvline(mean_gibbs - std_dev_gibbs, color='green', linestyle='dashed', linewidth=2, label
         plt.axvline(mean_gibbs + std_dev_gibbs, color='green', linestyle='dashed', linewidth=2)
         plt.axvline(x, color='black', linestyle='solid', linewidth=2, label=f'true value: {x:.2f}')
         plt.legend()
         plt.xlabel('probability of $p(c^+|f^+,g^+)$')
         plt.ylabel('Frequency')
         plt.title('Histogram of Gibbs sampling')
         plt.grid(True)
         plt.show()
```



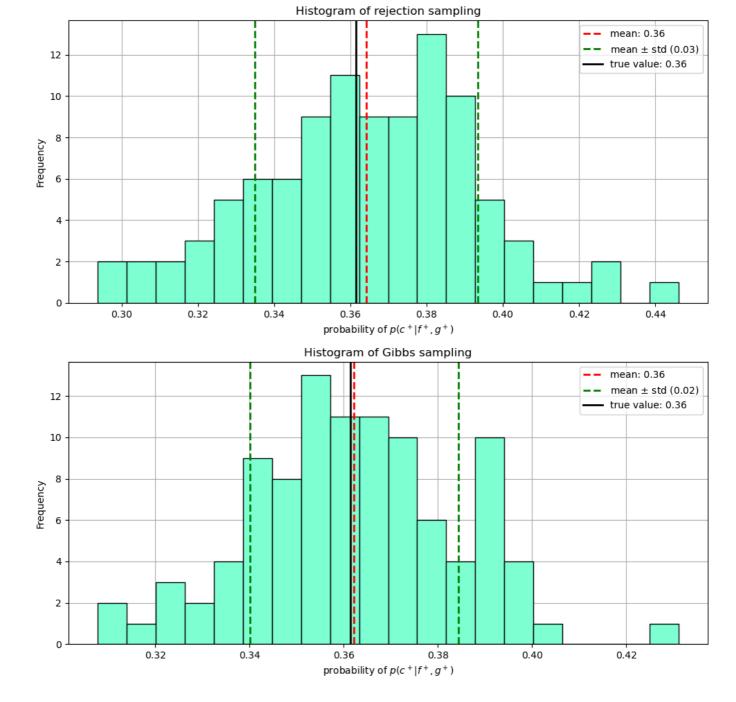
1.5. Poređenje metoda

Prikazaćemo rezultate obe metode jedno ispod drugog.

```
In [37]: import matplotlib.pyplot as plt

x = 0.361364
mean_rejection = np.mean(rejection_estimations)
std_dev_rejection = np.std(rejection_estimations)
mean_gibbs = np.mean(gibbs_estimations)
std_dev_gibbs = np.std(gibbs_estimations)
```

```
plt.figure(figsize=(10,10))
plt.subplot(211)
plt.rc('axes', axisbelow=True)
plt.hist(rejection_estimations, color='aquamarine', bins=20, edgecolor='black')
plt.axvline(mean_rejection, color='red', linestyle='dashed', linewidth=2, label=f'mean: {mean_
plt.axvline(mean_rejection - std_dev_rejection, color='green', linestyle='dashed', linewidth=
plt.axvline(mean_rejection + std_dev_rejection, color='green', linestyle='dashed', linewidth=
plt.axvline(x, color='black', linestyle='solid', linewidth=2, label=f'true value: {x:.2f}')
plt.xlabel('probability of p(c^+|f^+,g^+)')
plt.ylabel('Frequency')
plt.title('Histogram of rejection sampling')
plt.legend()
plt.grid(True)
plt.subplot(212)
plt.rc('axes', axisbelow=True)
plt.hist(gibbs_estimations, color='aquamarine', bins=20, edgecolor='black')
plt.axvline(mean_gibbs, color='red', linestyle='dashed', linewidth=2, label=f'mean: {mean_gib|
plt.axvline(mean_gibbs - std_dev_gibbs, color='green', linestyle='dashed', linewidth=2, label
plt.axvline(mean_gibbs + std_dev_gibbs, color='green', linestyle='dashed', linewidth=2)
plt.axvline(x, color='black', linestyle='solid', linewidth=2, label=f'true value: {x:.2f}')
plt.xlabel('probability of p(c^+|f^+,g^+)')
plt.ylabel('Frequency')
plt.title('Histogram of Gibbs sampling')
plt.legend()
plt.grid(True)
plt.tight_layout()
plt.show()
```



Zadatak 2: Čestični filtar

Čestični filtar predstavlja tehniku temporalnog probabilističkog rezonovanja. Ideja je da generisanjem više raznih čestica procenimo putanju kretanja objekta. Data su nam merenja rastojanja i ugla od merne stanice u fajlu 'observations.csv'.

2.1. Učitavanje .csv fajla

Učitaćemo podatke iz fajla. Merenja se vrše na svakih 1s i zašumljena su. Kako su podaci zašumljeni ćemo opisati u nastavku nakon što opišemo kako modelujemo stanja filtra.

```
In [9]: # Reading data from observations.csv
import numpy as np
from random import uniform, random, choice
from math import pi, cos, sin, sqrt, exp, atan
import matplotlib.pyplot as plt

data = np.loadtxt('./observations.csv', delimiter=',')
```

```
T,_ = data.shape
rho = data[:,0]
theta = data[:,1]

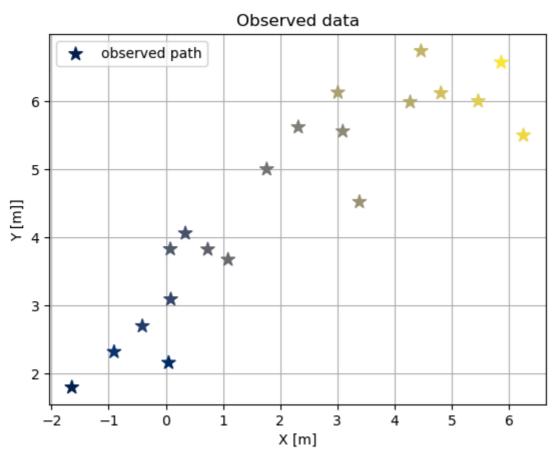
# Plotting the observed data
x = [rho[i]*cos(theta[i]) for i in range(T)]
y = [rho[i]*sin(theta[i]) for i in range(T)]

# Color gradient
colors = np.linspace(0, 1, T)

plt.scatter(x, y, c=colors, cmap='cividis', label='observed path', marker='*', s=100)

plt.title('Observed data')
plt.xlabel('X [m]')
plt.ylabel('Y [m]]')
plt.legend()
plt.grid('on')

plt.show()
```



2.2. Model česticnog filtra

Ovde treba opisati model za X_t^i i svaki korak koji sprovodimo, kao sto sam u svesci opisao

```
Stanje filtra: X_t = [x_t, y_t, \Omega_t, k_t]
```

x,y - pozicije cestice u XY ravni

 Ω - ugao vektora brzine u odnosu na x osu

k - broj sekundi od poslednjeg menjanja smera brzine

Formiraćemo N=100 cestica sa inicijalnom raspodelom (opisanu ispod) i u svakoj iteraciji cemo vrsiti sledece korake:

- ullet Predikcija: $X_{t|t-1}^{(i)} \sim p(X_t|X_{t-1}^{(i)})$
- Tezinjenje: $w_t^{(i)} \propto p(E_t|X_{t|t-1}^{(i)})$
- Estimacija polozaja: $E(X_t|E_{1:t})$
- Reuzorkovanje

2.3. Inicijalizacija čestičnog filtra: $X_0^{(i)} \sim p(X_0)$

```
egin{split} x_0^{(i)} &\sim Unif[-2,2] \ & \ y_0^{(i)} &\sim Unif[-2,2] \ & \ \Omega_0^{(i)} &\sim Unif[0,2\pi] \ & \ k_0^{(i)} &\sim \left(egin{array}{ccccc} 0 & 1 & 2 & 3 & 4 \ 1/5 & 1/5 & 1/5 & 1/5 & 1/5 \end{array}
ight) \end{split}
```

2.4. Predikcija $X_{1|0}^{(i)} \sim p(X_1^{(i)}|X_0^{(i)})$

```
\begin{split} x_{1|0}^{(i)} &= x_0^{(i)} + 0.5 \cdot cos(\omega_0^{(i)}) \\ y_{1|0}^{(i)} &= y_0^{(i)} + 0.5 \cdot sin(\omega_0^{(i)}) \\ p(\text{ promena } \Omega|k_0^{(i)}) &= 0.2 \cdot (k_0^{(i)} + 1) \\ if \text{ promena } \Omega: \\ &\Omega_{1|0}^{(i)} &= \Omega_0^{(i)} + \Delta, \Delta \sim Unif[-\pi/6, \pi/6], \Omega_{1|0}^{(i)} \in [0, 2\pi) \\ k_{1|0}^{(i)} &= 0 \\ else: \\ &\Omega_{1|0}^{(i)} &= \Omega_0^{(i)} \\ k_{1|0}^{(i)} &= k_0^{(i)} + 1 \end{split}
```

```
In [11]: predictions = np.empty((T, N, 4), dtype=float)

for i in range(N):
    x_0, y_0, omega_0, k_0 = particles[0, i]
```

```
x_1 = x_0 + 0.5*cos(omega_0)
y_1 = y_0 + 0.5*sin(omega_0)
r = random()
omega_1 = None
k_1 = None
if r < 0.2*(k_0+1):
    # Change direction of motion
    delta = uniform(-pi/6, pi/6)
    omega_1 = omega_0 + delta
    if omega_1>2*pi:
        omega_1-=2*pi
    if omega_1<0:</pre>
       omega 1+=2*pi
    k_1 = 0
else:
    omega_1 = omega_0
    k_1 = k_0 + 1
predictions[0, i] = [x_1, y_1, omega_1, k_1]
```

2.5. Tezinjenje prve iteracije, $w_1^{(i)} \propto p(E_1|X_{1|0}^{(i)})$

```
egin{aligned} E1 &= [
ho_1, 	heta_1] \ w_1^{(i)} &\propto p(
ho_1 | x_{1|0}^{(i)}, y_{1|0}^{(i)}, \Omega_{1|0}^{(i)}) \cdot p(	heta_1 | x_{1|0}^{(i)}, y_{1|0}^{(i)}) \ 
ho | x, y, \Omega &\sim \mathcal{N}(\sqrt{x^2 + y^2}, \sigma^2), \sigma = 0.3 \cdot (2 - sin(	heta - \Omega)) \ 	heta | x, y &\sim Laplace(arctan rac{y}{x}, b), 2b^2 = \left(rac{\pi}{36}
ight)^2 \end{aligned}
```

```
In [12]: weights = np.empty((T, N), dtype=float)
         rho 1 = rho[0]
         theta 1 = theta[0]
         b = pi/36/sqrt(2)
         for i in range(N):
             x_1, y_1, omega_1, k_1 = predictions[0, i]
             sigma = 0.3*(2-sin(theta 1-omega 1))
             aposteriori = 1
             aposteriori *= 1/sigma/sqrt(2*pi)
             temp = rho_1 - sqrt(x_1 ** 2 + y_1 ** 2)
             temp **= 2
             temp *= -1/2/(sigma ** 2)
             aposteriori *= exp(temp)
             aposteriori *= 1/2/b
             temp = theta_1-np.arctan2(y_1, x_1)
             temp = -abs(temp)/b
             aposteriori *= exp(temp)
             weights[0, i] = aposteriori
         weights[0] = weights[0]/np.sum(weights[0])
```

```
In [13]: # Prvi neophodni grafici: Tezine cestica posle prve iter
import matplotlib.pyplot as plt

plt.figure(figsize=(20,6))
```

```
plt.subplot(121)
  plt.scatter(particles[0,:,0], particles[0,:,1], color='blue', label='Data points')
  plt.scatter([rho_1*cos(theta_1)],[rho_1*sin(theta_1)], color='red', marker='x')
  plt.title('Initialised particle positions')
  plt.xlabel('X [m]')
  plt.ylabel('Y [m]]')
 plt.grid('on')
 plt.subplot(122)
 current_weights = weights[0]
 max_weight = max(current_weights)
 normalized_weights = [weight / max_weight * 100 for weight in current_weights]
 plt.scatter(predictions[0,:,0], predictions[0,:,1], s=normalized_weights, color='blue', label
 plt.scatter([rho_1*cos(theta_1)],[rho_1*sin(theta_1)], color='red', marker='x')
 e_x = 0
 e_y = 0
 for i in range(N):
      e_x += weights[0,i]*predictions[0,i,0]
      e_y += weights[0,i]*predictions[0,i,1]
  plt.scatter([e_x], [e_y],marker='*', color='black')
 plt.title('Particle positions at t=1')
 plt.xlabel('X [m]')
 plt.ylabel('Y [m]]')
 plt.grid('on')
 plt.show()
                                                                       Particle positions at t=1
                                                               ×
                                                    <u>E</u> 0
0.0
 -0.5
```

2.6. Druga iteracija bez reuzorkovanja

2.6.1. Nereuzorkovanje

-1.5

```
In [14]: predictions_no_resample = np.empty((N,4),dtype=float)
weights_no_resample = np.empty(N,dtype=float)
```

2.6.2 Predikcija $X_{2|1}^{(i)} \sim p(X_2^{(i)}|X_1^{(i)})$

```
omega_2 = omega_1 + delta
if omega_2>2*pi:
    omega_2-=2*pi
if omega_2<0:
    omega_2+=2*pi
    k_2 = 0
else:
    omega_2 = omega_1
    k_2 = k_1 + 1

predictions_no_resample[i] = [x_2, y_2, omega_2, k_2]</pre>
```

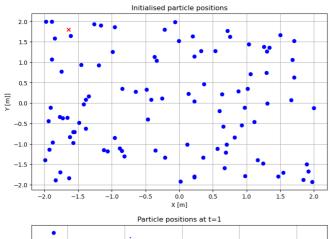
2.6.3. Tezinjenje druge iteracije, $w_2^{(i)} \propto p(E_2|X_{2|1}^{(i)})$

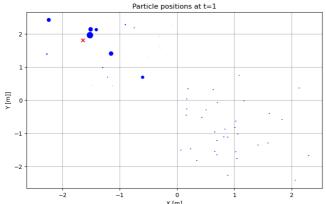
```
In [16]: rho_2 = rho[1]
         theta_2 = theta[1]
         b = pi/36/sqrt(2)
         for i in range(N):
             x_2, y_2, omega_2, k_2 = predictions_no_resample[i]
             sigma = 0.3*(2-sin(theta_2-omega_2))
             aposteriori = 1
             aposteriori *= 1/sigma/sqrt(2*pi)
             temp = rho_2 - sqrt(x_2 ** 2 + y_2 ** 2)
             temp **= 2
             temp *= -1/2/(sigma ** 2)
             aposteriori *= exp(temp)
             aposteriori *= 1/2/b
             temp = theta_2-np.arctan2(y_2, x_2)
             temp = -abs(temp)/b
             aposteriori *= exp(temp)
             weights_no_resample[i] = aposteriori
         weights_no_resample = weights_no_resample/sum(weights_no_resample)
```

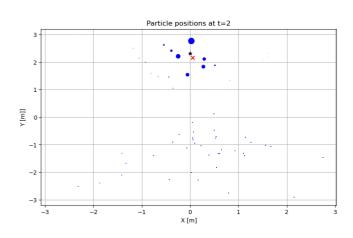
```
In [17]:
         # Drugi neophodni grafici: Tezine cestica posle druge iter bez reuzorkovanja
         import matplotlib.pyplot as plt
         plt.figure(figsize=(20,12))
         plt.subplot(221)
         plt.scatter(particles[0,:,0],\ particles[0,:,1],\ color='blue',\ label='Data\ points')
         plt.scatter([rho_1*cos(theta_1)],[rho_1*sin(theta_1)], color='red', marker='x')
         plt.title('Initialised particle positions')
         plt.xlabel('X [m]')
         plt.ylabel('Y [m]]')
         plt.grid('on')
         plt.subplot(223)
         current_weights = weights[0]
         max_weight = max(current_weights)
         normalized_weights = [weight / max_weight * 100 for weight in current_weights]
         plt.scatter(predictions[0,:,0], predictions[0,:,1], s=normalized_weights, color='blue', label
         plt.scatter([rho_1*cos(theta_1)],[rho_1*sin(theta_1)], color='red', marker='x')
         plt.title('Particle positions at t=1')
         plt.xlabel('X [m]')
         plt.ylabel('Y [m]]')
         plt.grid('on')
         e x = 0
         e_y = 0
```

```
for i in range(N):
    e_x += weights_no_resample[i]*predictions_no_resample[i,0]
    e_y += weights_no_resample[i]*predictions_no_resample[i,1]

plt.subplot(224)
current_weights = weights_no_resample
max_weight = max(current_weights)
normalized_weights = [weight / max_weight * 100 for weight in current_weights]
plt.scatter(predictions_no_resample[:,0], predictions_no_resample[:,1], s=normalized_weights,
plt.scatter([rho_2*cos(theta_2)],[rho_2*sin(theta_2)], color='red', marker='x')
plt.scatter([e_x],[e_y],marker='*',color='black')
plt.title('Particle positions at t=2')
plt.xlabel('X [m]')
plt.ylabel('Y [m]]')
plt.grid('on')
```







2.7. Druga iteracija sa reuzorkovanjem

$$j \sim \left(egin{array}{cccc} 0 & 1 & \dots & N-1 \ w_t^{(0)} & w_t^{(1)} & \dots & w_t^{(N)} \end{array}
ight)$$

2.7.1. Reuzorkovanje

```
In [18]: # Koristimo nizove weights[0], indexes i running_sum

# Pravimo listu takvu da...
# Unsorted list of floats
unsorted_list = weights[0]

# Pair each element with its index using enumerate
indexed_list = list(enumerate(unsorted_list))

# Sort the list of pairs based on the values (second element of the pairs)
sorted_indexed_list = sorted(indexed_list, key=lambda x: x[1])
```

```
# Extract the indices from the sorted pairs
indexes = [index for index, value in sorted_indexed_list]
sorted_weights = [value for index, value in sorted_indexed_list]

running_sum = np.cumsum(sorted_weights)

for i in range(N):
    r = random()
    j = 0
    while(running_sum[j]<r):
        j+=1
    particles[1,i] = predictions[0, indexes[j]]</pre>
```

2.7.2. Predikcija $X_{2|1}^{(i)} \sim p(X_2^{(i)}|X_1^{(i)})$

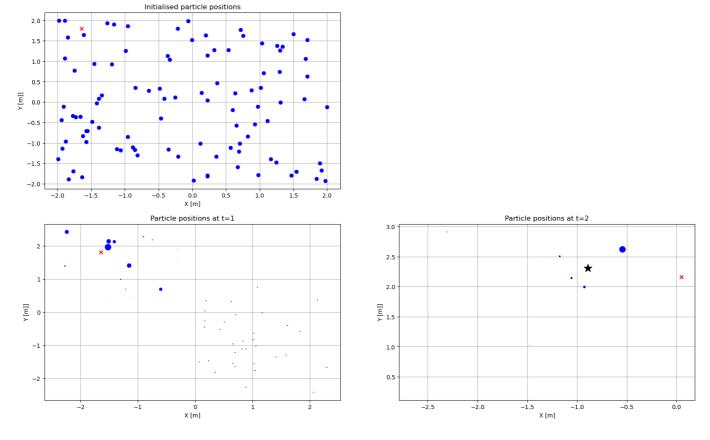
```
In [19]: for i in range(N):
             x_1, y_1, omega_1, k_1 = particles[1, i]
             x_2 = x_1 + 0.5*cos(omega_1)
             y_2 = y_1 + 0.5*sin(omega_1)
             r = random()
             omega_2 = None
             k_2 = None
              if r < 0.2*(k_1+1):
                  # Change direction of motion
                  delta = uniform(-pi/6, pi/6)
                  omega_2 = omega_1 + delta
                  if omega_2>2*pi:
                      omega_2-=2*pi
                  if omega_2<0:</pre>
                      omega_2+=2*pi
                  k_2 = 0
              else:
                  omega_2 = omega_1
                  k_2 = k_1 + 1
              predictions[1, i] = [x_2, y_2, omega_2, k_2]
```

2.7.3. Tezinjenje druge iteracije, $w_2^{(i)} \propto p(E_2|X_{2|1}^{(i)})$

```
In [20]:
         rho_2 = rho[1]
         theta 2 = theta[1]
         b = pi/36/sqrt(2)
         for i in range(N):
             x_2, y_2, omega_2, k_2 = predictions[1, i]
             sigma = 0.3*(2-sin(theta_2-omega_2))
             aposteriori = 1
             aposteriori *= 1/sigma/sqrt(2*pi)
             temp = rho_2 - sqrt(x_2 ** 2 + y_2 ** 2)
             temp **= 2
             temp *= -1/2/(sigma ** 2)
             aposteriori *= exp(temp)
             aposteriori *= 1/2/b
             temp = theta_2-np.arctan2(y_2, x_2)
             temp = -abs(temp)/b
             aposteriori *= exp(temp)
             weights[1, i] = aposteriori
```

```
weights[1] = weights[1]/np.sum(weights[1])
```

```
In [21]:
         # Drugi neophodni grafici: Tezine cestica posle druge iter sa reuzorkovanjem
         import matplotlib.pyplot as plt
         plt.figure(figsize=(20,12))
         plt.subplot(221)
         plt.scatter(particles[0,:,0], particles[0,:,1], color='blue', label='Data points')
         plt.scatter([rho_1*cos(theta_1)],[rho_1*sin(theta_1)], color='red', marker='x')
         plt.title('Initialised particle positions')
         plt.xlabel('X [m]')
         plt.ylabel('Y [m]]')
         plt.grid('on')
         plt.subplot(223)
         current_weights = weights[0]
         max_weight = max(current_weights)
         normalized_weights = [weight / max_weight * 100 for weight in current_weights]
         plt.scatter(predictions[0,:,0], predictions[0,:,1], s=normalized_weights, color='blue', label
         plt.scatter([rho_1*cos(theta_1)],[rho_1*sin(theta_1)], color='red', marker='x')
         plt.title('Particle positions at t=1')
         plt.xlabel('X [m]')
         plt.ylabel('Y [m]]')
         plt.grid('on')
         e_x = 0
         e_y = 0
         for i in range(N):
             e_x += weights[1,i]*predictions[1,i,0]
             e_y += weights[1,i]*predictions[1,i,1]
         plt.subplot(224)
         current_weights = weights[1]
         max_weight = max(current_weights)
         normalized_weights = [weight / max_weight * 100 for weight in current_weights]
         plt.scatter(predictions[1,:,0], predictions[1,:,1], s=normalized_weights, color='blue', label
         plt.scatter([rho_2*cos(theta_2)],[rho_2*sin(theta_2)], color='red', marker='x')
         plt.scatter([e_x],[e_y],color='black',marker='*',s=200)
         plt.title('Particle positions at t=2')
         plt.xlabel('X [m]')
         plt.ylabel('Y [m]]')
         plt.grid('on')
         plt.show()
```



Komentar: Uočava se efekat 'osiromašenja' uzorka (mana postupka reuzorkovanja).

2.8. Izvršavanje čestičnog filtra

Izvršene su prve 2 iteracije, potrebno je uraditi jos T-2

```
In [22]:
         for t in range(2,T):
              # Reuzorkovanje
              unsorted_list = weights[t-1]
              indexed list = list(enumerate(unsorted list))
              sorted_indexed_list = sorted(indexed_list, key=lambda x: x[1])
              indexes = [index for index, value in sorted_indexed_list]
              sorted_weights = [value for index, value in sorted_indexed_list]
              running_sum = np.cumsum(sorted_weights)
              for i in range(N):
                  r = random()
                  j = 0
                  while(running_sum[j]<r):</pre>
                  particles[t, i] = predictions[t-1, indexes[j]]
              # Predikcija
              for i in range(N):
                  x_t, y_t, omega_t, k_t = particles[t, i]
                  x_pred = x_t + 0.5*cos(omega_t)
                  y_pred = y_t + 0.5*sin(omega_t)
                  r = random()
                  omega_pred = None
                  k_pred = None
                  if r < 0.2*(k_t+1):
                      # Change direction of motion
                      delta = uniform(-pi/6, pi/6)
                      omega_pred = omega_t + delta
                      if omega_pred>2*pi:
                          omega_pred-=2*pi
```

```
if omega_pred<0:</pre>
            omega_pred+=2*pi
        k_pred = 0
    else:
        omega_pred = omega_t
        k_pred = k_t + 1
    predictions[t, i] = [x_pred, y_pred, omega_pred, k_pred]
# Tezinjenje
rho_t = rho[t]
theta_t = theta[t]
b = pi/36/sqrt(2)
for i in range(N):
    x_pred, y_pred, omega_pred, k_pred = predictions[t, i]
    sigma = 0.3*(2-sin(theta_t-omega_pred))
   w = 1
   w *= 1/sigma/sqrt(2*pi)
   temp = rho_t - sqrt(x_pred ** 2 + y_pred ** 2)
   temp **= 2
   temp *= -1/2/(sigma ** 2)
   w *= exp(temp)
   w *= 1/2/b
   temp = theta_t-np.arctan2(y_pred, x_pred)
    temp = -abs(temp)/b
    w *= exp(temp)
    weights[t, i] = w
weights[t] = weights[t]/np.sum(weights[t])
```

2.9. Estimacije $E(X_t | E_{1:t})$

$$egin{align} E(x_t|E_{1:t}) &pprox \sum\limits_{i=1}^N w^{(i)} x_t^{(i)} \ &E(y_t|E_{1:t}) pprox \sum\limits_{i=1}^N w^{(i)} y_t^{(i)} \ &E(
ho_t|E_{1:t}) = \sqrt{E_{x_t}^2 + E_{y_t}^2} \ &E(heta_t|E_{1:t}) = arctanrac{E_{y_t}}{E_{x_t}} \ \end{array}$$

```
In [23]: x_estimated = np.empty(T)
y_estimated = np.empty(T)
uncertainty_sigma = np.empty(T, dtype=float)

for t in range(T):
    e_x = 0
    e_y = 0
    e_omega = 0
    for i in range(N):
        e_x += weights[t,i]*predictions[t,i,0]
        e_y += weights[t,i]*predictions[t,i,1]
        e_omega += predictions[t,i,2]
    e_omega /= N

uncertainty_sigma[t] = 0.3*(2-sin(theta[t]-e_omega))
```

```
y_estimated[t] = e_y

In [24]: plt.figure(figsize=(8,6))
    plt.scatter(x, y, color='blue', marker='.', label='observed path')

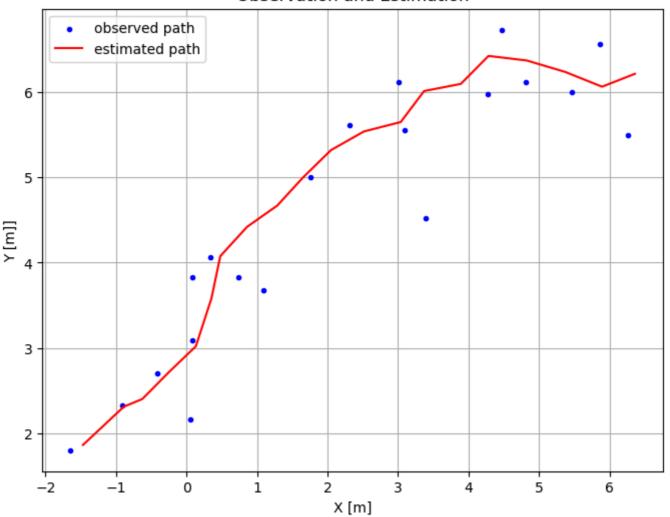
plt.plot(x_estimated, y_estimated, color='red', label='estimated path')

plt.title('Observation and Estimation')
    plt.xlabel('X [m]')
    plt.ylabel('Y [m]]')
    plt.grid('on')
    plt.legend()

plt.show()
```

 $x_{estimated[t]} = e_x$

Observation and Estimation



```
In [25]: rho_estimated = [sqrt(x_estimated[t]**2+y_estimated[t]**2) for t in range(T)]

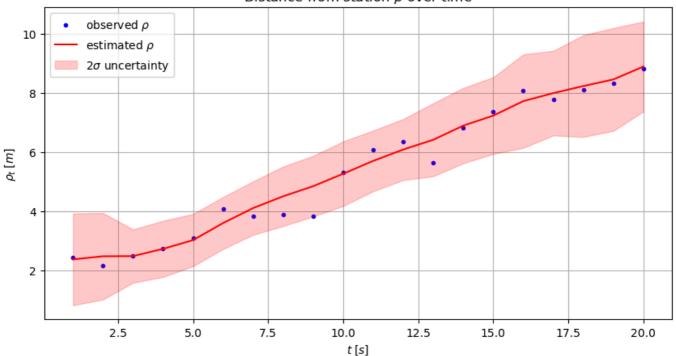
plt.figure(figsize=(10,5))
    t = range(1,21)
    plt.scatter(t, rho, marker='.', color='blue', label=r'observed $\rho$')
    plt.plot(t, rho_estimated, color='red', label=r'estimated $\rho$')

# 2-sigma uncertainty
    variance_upper = rho_estimated+2*uncertainty_sigma
    variance_lower = rho_estimated-2*uncertainty_sigma
    # plt.plot(t, variance_upper, linewidth=0.5)
    # plt.plot(t, variance_lower, linewidth=0.5)
    plt.fill_between(t,variance_lower,variance_upper,color='red',alpha=0.2, label=r'$2\sigma$ unco

plt.xlabel(r'$t \; [s]$')
    plt.ylabel(r'$\rho_t \; [m]$')
    plt.title(r'Distance from station $\rho$ over time')
```

```
plt.grid('on')
plt.legend()
plt.show()
```

Distance from station ρ over time



```
In [26]: theta_estimated = np.array([np.arctan2(y_estimated[t],x_estimated[t]) for t in range(T)])

plt.figure(figsize=(10,5))
t = range(1,21)
plt.scatter(t, theta, marker='.', color='blue', label=r'observed $\theta$')
plt.plot(t, theta_estimated, color='red', label=r'estimated $\theta$')
# 2-sigma uncertainty
variance_upper = theta_estimated+2*pi/36
variance_lower = theta_estimated-2*pi/36
plt.fill_between(t,variance_lower,variance_upper, color='red',alpha=0.2, label=r'$2\sigma$ un

plt.xlabel(r'$t \;[s]$')
plt.ylabel(r'$\theta_t \; [rad]$')
plt.title(r'Angle from station $\theta$ over time')
plt.grid('on')
plt.legend()

plt.show()
```

